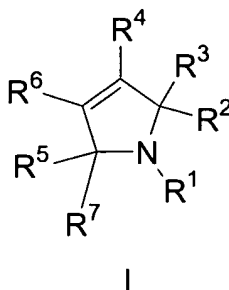


In the claims:

1. (Currently amended) A compound of Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof,

wherein:

- a is 0 or 1;
b is 0 or 1;
m is 0, 1, or 2;
n is 0;
r is 0 or 1;
s is 0 or 1;

R¹ is selected from:

- 1) (C₁-C₆-alkylene)_n(C=O)C₁-C₁₀ alkyl, and
- 2) —(C₁-C₆-alkylene)_n(C=O)aryl,
- 3) —(C₁-C₆-alkylene)_n(C=O)C₂-C₁₀-alkenyl,
- 4) —(C₁-C₆-alkylene)_n(C=O)C₂-C₁₀-alkynyl,
- 5) —(C₁-C₆-alkylene)_n(C=O)C₃-C₈-cycloalkyl,
- 6) —(C₁-C₆-alkylene)_n(C=O)heterocyclyl,
- 7) (C₁-C₆-alkylene)_n(C=O)NR^cR^{c'},
- 8) —(C₁-C₆-alkylene)_nSO₂NR^eR^{e'},
- 9) —(C₁-C₆-alkylene)_nSO₂C₁-C₁₀-alkyl,
- 10) —(C₁-C₆-alkylene)_nSO₂-aryl,
- 11) —(C₁-C₆-alkylene)_nSO₂-heterocyclyl,

- 12) $(C_1-C_6\text{-alkylene})_nSO_2-C_3-C_8\text{-cycloalkyl}$;
- 13) $(C_1-C_6\text{-alkylene})_nP(=O)R^dR^{d'}$;
- 14) aryl ;
- 15) heterocycetyl ;
- 16) $C_1-C_{10}\text{-alkyl}$;
- 17) $(C_1-C_6\text{-alkylene})_n(C=O)O-C_1-C_{10}\text{-alkyl}$;
- 18) $(C_1-C_6\text{-alkylene})_n(C=O)O\text{-aryl}$;
- 19) $(C_1-C_6\text{-alkylene})_n(C=O)O-C_2-C_{10}\text{-alkenyl}$;
- 20) $(C_1-C_6\text{-alkylene})_n(C=O)O-C_2-C_{10}\text{-alkynyl}$;
- 21) $(C_1-C_6\text{-alkylene})_n(C=O)O-C_3-C_8\text{-cycloalkyl}$;
- 22) $(C_1-C_6\text{-alkylene})_n(C=O)O\text{-heterocycetyl}$;

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocycetyl is optionally substituted with one or more substituents selected from R^{10} ;

R^2 and R^6 are independently selected from:

- 1) aryl ,
- 2) $C_1-C_6\text{-aralkyl}$,
- 3) $C_3-C_8\text{-cycloalkyl}$, and
- 4) heterocycetyl ;

said aryl, cycloalkyl, aralkyl and heterocycetyl is optionally substituted with one or more substituents selected from R^{10} ;

R^3 is selected from:

- 1) $C_1-C_{10}\text{-alkyl-O-R}^g$,
- 2) $C_2-C_{10}\text{-alkenyl-O-R}^g$,
- 3) $C_2-C_{10}\text{-alkynyl-O-R}^g$,
- 4) $(C_1-C_6\text{-alkylene})_n C_3-C_8\text{-cycloalkyl-O-R}^g$,
- 5) $C_1-C_{10}\text{-alkyl-(C=O)}_b\text{-NR}^fR^{f'}$,
- 6) $C_2-C_{10}\text{-alkenyl-(C=O)}_b\text{-NR}^fR^{f'}$,
- 7) $C_2-C_{10}\text{-alkynyl-(C=O)}_b\text{-NR}^fR^{f'}$,
- 8) $(C_1-C_6\text{-alkylene})_n C_3-C_8\text{-cycloalkyl-(C=O)}_b\text{-NR}^fR^{f'}$,
- 9) $C_1-C_{10}\text{-alkyl-S(O)}_m\text{-R}^g$,
- 10) $C_2-C_{10}\text{-alkenyl-S(O)}_m\text{-R}^g$,

11) ~~C₂-C₁₀ alkynyl S(O)_m R⁸,~~

12) ~~(C₁-C₆ alkylene)_n C₃-C₈ cycloalkyl S(O)_m R⁸,~~

said alkyl, ~~alkenyl, alkynyl and cycloalkyl~~ are is optionally substituted with one or more substituents selected from R¹⁰;

R⁴ is selected from:

- 1) H, and
- 2) C₁-C₁₀ alkyl,
- 3) ~~aryl,~~
- 4) ~~C₂-C₁₀ alkenyl,~~
- 5) ~~C₂-C₁₀ alkynyl,~~
- ~~6) C₁-C₆ perfluoroalkyl,~~
- ~~7) C₁-C₆ aralkyl,~~
- 8) ~~C₃-C₈ cycloalkyl, and~~
- ~~9) heterocyclyl,~~

said alkyl, ~~aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R¹⁰;

R⁵ and R⁷ are independently selected from:

- 1) H, and
- 2) C₁-C₁₀ alkyl,
- 3) ~~aryl,~~
- 4) ~~C₂-C₁₀ alkenyl,~~
- 5) ~~C₂-C₁₀ alkynyl,~~
- ~~6) C₁-C₆ perfluoroalkyl,~~
- ~~7) C₁-C₆ aralkyl,~~
- 8) ~~C₃-C₈ cycloalkyl, and~~
- ~~9) heterocyclyl,~~

said alkyl, ~~aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R¹⁰; or

~~R⁵ and R⁷ are combined to form an oxo or a sulfoxo;~~

R¹⁰ is independently selected from:

- 1) (C=O)_aO_bC₁-C₁₀ alkyl,
- 2) (C=O)_aO_baryl,
- 3) C₂-C₁₀ alkenyl,
- 4) C₂-C₁₀ alkynyl,
- 5) (C=O)_aO_b heterocyclyl,
- 6) CO₂H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O_bC₁-C₆ perfluoroalkyl,
- 11) O_a(C=O)_bNR¹²R¹³,
- 12) S(O)_mR^a,
- 13) S(O)₂NR¹²R¹³,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R¹²R¹³, or
- 17) (C=O)_aO_bC₃-C₈ cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R¹¹;

R¹¹ is selected from:

- 1) (C=O)_rO_s(C₁-C₁₀)alkyl,
- 2) O_r(C₁-C₃)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C₂-C₁₀)alkenyl,
- 8) (C₂-C₁₀)alkynyl,
- 9) (C=O)_rO_s(C₃-C₆)cycloalkyl,
- 10) (C=O)_rO_s(C₀-C₆)alkylene-aryl,
- 11) (C=O)_rO_s(C₀-C₆)alkylene-heterocyclyl,
- 12) (C=O)_rO_s(C₀-C₆)alkylene-N(R^b)₂,

- 13) $C(O)R^a$,
- 14) $(C_0-C_6)\text{alkylene-CO}_2R^a$,
- 15) $C(O)H$,
- 16) $(C_0-C_6)\text{alkylene-CO}_2H$, and
- 17) $C(O)N(R^b)_2$,
- 18) $S(O)_mR^a$, and
- 19) $S(O)_2N(R^b)_2$;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R^b , OH, $(C_1-C_6)\text{alkoxy}$, halogen, CO_2H , CN, $O(C=O)C_1-C_6$ alkyl, oxo, NO_2 and $N(R^b)_2$;

R^{12} and R^{13} are independently selected from:

- 1) H,
- 2) $(C=O)O_bC_1-C_{10}$ alkyl,
- 3) $(C=O)O_bC_3-C_8$ cycloalkyl,
- 4) $(C=O)O_b\text{aryl}$,
- 5) $(C=O)O_b\text{heterocyclyl}$,
- 6) C_1-C_{10} alkyl,
- 7) aryl,
- 8) C_2-C_{10} alkenyl,
- 9) C_2-C_{10} alkynyl,
- 10) heterocyclyl,
- 11) C_3-C_8 cycloalkyl,
- 12) SO_2R^a , and
- 13) $(C=O)NR^b_2$,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R^{11} , or

R^{12} and R^{13} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R^{11} ;

R^a is independently selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^b is independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fR^{f'} or S(O)₂R^a, optionally substituted with one, two or three substituents selected from R¹¹;

R^c and R^{c'} are independently selected from: H, (C₁-C₆)alkyl, ~~aryl~~, heterocyclyl and (C₃-C₆)cycloalkyl, optionally substituted with one, two or three substituents selected from R¹¹; or

R^c and R^{c'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^d and R^{d'} are independently selected from: (C₁-C₆)alkyl, (C₁-C₆)alkoxy and NR^b₂, or

R^d and R^{d'} can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 3-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR^e, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹; and

R^e is selected from: H and (C₁-C₆)alkyl, optionally substituted with one, two or three substituents selected from R¹¹;

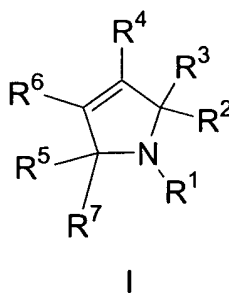
R^f and R^{f'} are independently selected from: H, (C₁-C₆)alkyl, aryl, NH₂, OH, OR^a, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fR^{f'}, S(O)₂R^a and -(C₁-C₆)alkyl-N(R^b)₂, wherein the alkyl is optionally substituted with one, two or three substituents selected from R¹¹; or

R^f and R^{f'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R^{11} ;

R^8 is selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂.

2. (Currently amended) The compound according to Claim 1 of Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof,

wherein:

a is 0 or 1;
b is 0 or 1;
m is 0, 1, or 2;
n is 0 or 1;
r is 0 or 1;
s is 0 or 1;

R^1 is selected from:

- 1) (C₁-C₆-alkylene)_n(C=O)C₁-C₁₀ alkyl,
- 2) (C₁-C₆-alkylene)_n(C=O)aryl,
- 3) (C₁-C₆-alkylene)_n(C=O)C₂-C₁₀ alkenyl,
- 4) (C₁-C₆-alkylene)_n(C=O)C₂-C₁₀ alkynyl,
- 5) (C₁-C₆-alkylene)_n(C=O)C₃-C₈ cycloalkyl,
- 6) (C₁-C₆-alkylene)_n(C=O)heterocyclyl,
- 2) 7) (C₁-C₆-alkylene)_n(C=O)NR^cR^{c'},

- 7) —(C1-C6-alkylene)_nSO₂NReRe',
- 8) —(C1-C6-alkylene)_nSO₂C1-C10-alkyl,
- 9) —(C1-C6-alkylene)_nSO₂-aryl,
- 10) —(C1-C6-alkylene)_nSO₂-heterocycetyl,
- 11) —(C1-C6-alkylene)_nSO₂-C3-C8-cycloalkyl,
- 12) —(C1-C6-alkylene)_nP(=O)RdRd',
- 13) —aryl,
- 14) —heterocycetyl,
- 15) —C1-C10-alkyl,
- 16) —(C1-C6-alkylene)_n(C=O)O-C1-C10-alkyl,
- 17) —(C1-C6-alkylene)_n(C=O)O-aryl,
- 18) —(C1-C6-alkylene)_n(C=O)O-C2-C10-alkenyl,
- 19) —(C1-C6-alkylene)_n(C=O)O-C2-C10-alkynyl,
- 20) —(C1-C6-alkylene)_n(C=O)O-C3-C8-cycloalkyl,
- 21) —(C1-C6-alkylene)_n(C=O)O-heterocycetyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocycetyl is optionally substituted with one or more substituents selected from R¹⁰;

R² and R⁶ are independently selected from:

- 1) aryl,
- 2) —C₁-C₆-aralkyl,
- 3) —C₃-C₈-cycloalkyl, and
- 4) —heterocycetyl,

said aryl, cycloalkyl, aralkyl and heterocycetyl is optionally substituted with one or more substituents selected from R¹⁰;

R³ is selected from:

- 1) C₁-C₁₀ alkyl-O-Rg,
- 2) —C₂-C₁₀ alkenyl-O-Rg,
- 3) —C₂-C₁₀ alkynyl-O-Rg,
- 4) —(C₁-C₆-alkylene)_nC₃-C₈-cycloalkyl-O-Rg,
- 2.5) C₁-C₁₀ alkyl-(C=O)_b-NR^fR^{f'},
- 6) —C₂-C₁₀ alkenyl-(C=O)_bNR^fR^{f'},
- 7) —C₂-C₁₀ alkynyl-(C=O)_bNR^fR^{f'},

8) —(C₁-C₆-alkylene)_nC₃-C₈-cycloalkyl (C=O)_bNR^fR^{f'};

9) —C₁-C₁₀-alkyl-S(O)_m-R^g;

10) —C₂-C₁₀-alkenyl-S(O)_m-R^g;

11) —C₂-C₁₀-alkynyl-S(O)_m-R^g;

12) —(C₁-C₆-alkylene)_nC₃-C₈-cycloalkyl-S(O)_m-R^g;

said alkyl, alkenyl, alkynyl and cycloalkyl are is optionally substituted with one or more substituents selected from R¹⁰;

R⁴ is selected from:

- 1) H,
- 2) C₁-C₁₀ alkyl,
- 3) —aryl,
- 4) —C₂-C₁₀-alkenyl,
- 5) —C₂-C₁₀-alkynyl,
- 6) —C₁-C₆-perfluoroalkyl,
- 7) —C₁-C₆-aralkyl,
- 8) —C₃-C₈-cycloalkyl, and

—9) —heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R⁵ and R⁷ are independently selected from:

- 1) H,
- 2) C₁-C₁₀ alkyl,
- 3) —aryl,
- 4) —C₂-C₁₀-alkenyl,
- 5) —C₂-C₁₀-alkynyl,
- 6) —C₁-C₆-perfluoroalkyl,
- 7) —C₁-C₆-aralkyl,
- 8) —C₃-C₈-cycloalkyl, and

—9) —heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰; or

~~R⁵ and R⁷ are combined to form an oxo or a sulfoxo;~~

R¹⁰ is independently selected from:

- 1) (C=O)_aO_bC₁-C₁₀ alkyl,
- 2) (C=O)_aO_baryl,
- 3) C₂-C₁₀ alkenyl,
- 4) C₂-C₁₀ alkynyl,
- 5) (C=O)_aO_b heterocyclyl,
- 6) CO₂H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O_bC₁-C₆ perfluoroalkyl,
- 11) O_a(C=O)_bNR¹²R¹³,
- 12) S(O)_mR^a,
- 13) S(O)₂NR¹²R¹³,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R¹²R¹³, or
- 17) (C=O)_aO_bC₃-C₈ cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R¹¹;

R¹¹ is selected from:

- 1) (C=O)_rO_s(C₁-C₁₀)alkyl,
- 2) O_r(C₁-C₃)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C₂-C₁₀)alkenyl,
- 8) (C₂-C₁₀)alkynyl,
- 9) (C=O)_rO_s(C₃-C₆)cycloalkyl,

- 10) $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-aryl}$,
- 11) $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-heterocyclyl}$,
- 12) $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-N(R}^b)_2$,
- 13) C(O)R^a ,
- 14) $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{R}^a$,
- 15) C(O)H ,
- 16) $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{H}$, and
- 17) $\text{C(O)N(R}^b)_2$,
- 18) $\text{S(O)}_m\text{R}^a$, and
- 19) $\text{S(O)}_2\text{N(R}^b)_2$;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R^b , OH, $(\text{C}_1\text{-C}_6)\text{alkoxy}$, halogen, CO_2H , CN, $\text{O(C=O)C}_1\text{-C}_6$ alkyl, oxo, and $\text{N(R}^b)_2$;

R^{12} and R^{13} are independently selected from:

- 1) H,
- 2) $(\text{C}=\text{O})\text{O}_b\text{C}_1\text{-C}_{10}$ alkyl,
- 3) $(\text{C}=\text{O})\text{O}_b\text{C}_3\text{-C}_8$ cycloalkyl,
- 4) $(\text{C}=\text{O})\text{O}_b\text{aryl}$,
- 5) $(\text{C}=\text{O})\text{O}_b\text{heterocyclyl}$,
- 6) $\text{C}_1\text{-C}_{10}$ alkyl,
- 7) aryl,
- 8) $\text{C}_2\text{-C}_{10}$ alkenyl,
- 9) $\text{C}_2\text{-C}_{10}$ alkynyl,
- 10) heterocyclyl,
- 11) $\text{C}_3\text{-C}_8$ cycloalkyl,
- 12) SO_2R^a , and
- 13) $(\text{C}=\text{O})\text{NR}^b_2$,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R^{11} , or

R^{12} and R^{13} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in

addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^a is independently selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^b is independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fR^{f'} or S(O)₂R^a, optionally substituted with one, two or three substituents selected from R¹¹;

R^c and R^{c'} are independently selected from: H, (C₁-C₆)alkyl, ~~aryl~~, heterocyclyl and (C₃-C₆)cycloalkyl, optionally substituted with one, two or three substituents selected from R¹¹; or

R^c and R^{c'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^d and R^{d'} are independently selected from: (C₁-C₆)alkyl, (C₁-C₆)alkoxy and NR^b₂, or

R^d and R^{d'} can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 3-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR^e, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹; and

R^e is selected from: H and (C₁-C₆)alkyl, optionally substituted with one, two or three substituents selected from R¹¹;

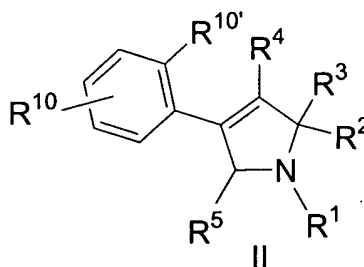
R^f and R^{f'} are independently selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂, or

R^f and R^{f'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R^{11} ;

R_8 is selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂.

3. (Currently amended) The compound according to Claim 2 of Formula II:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;
b is 0 or 1;
m is 0, 1, or 2;
r is 0 or 1;
s is 0 or 1;

R^1 is selected from:

- 1) (C=O)C₁-C₁₀ alkyl, and
- 2) —(C=O)aryl,
- 3) —(C=O)C₂-C₁₀ alkenyl,
- 4) —(C=O)C₂-C₁₀ alkynyl,
- 5) —(C=O)C₃-C₈ cycloalkyl,
- 6) —(C=O)heterocyclyl,
- 7) (C=O)NR^cR^{c'},
- 8) —SO₂NR^eR^{e'},
- 9) —SO₂C₁-C₁₀ alkyl,
- 10) —SO₂ aryl,
- 11) —SO₂ heterocyclyl,

~~12) —SO₂-C₃-C₈-cycloalkyl, and~~

~~13) —P(=O)R^dR^{d'},~~

said alkyl, ~~aryl, alkenyl, alkynyl, cycloalkyl, heteroaryl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R¹⁰;

R² is selected from:

1) aryl,

~~2) —C₁-C₆-aralkyl,~~

~~3) —C₃-C₈-cycloalkyl, and~~

~~4) —heterocyclyl,~~

said aryl, ~~cycloalkyl, aralkyl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R¹⁰;

R³ is selected from:

1) C₁-C₁₀ alkyl-O-R^g,

~~2) —C₃-C₈-cycloalkyl-O-R^g,~~

~~2 3) C₁-C₁₀ alkyl-NR^fR^{f'},~~

~~4) —C₃-C₈-cycloalkyl-NR^fR^{f'},~~

said alkyl ~~and cycloalkyl~~ are is optionally substituted with one or more substituents selected from R¹⁰;

R⁴ and R⁵ are independently selected from:

1) H, and

2) C₁-C₁₀ alkyl,

~~3) —aryl,~~

~~4) —C₂-C₁₀-alkenyl,~~

~~5) —C₂-C₁₀-alkynyl,~~

~~6) —C₁-C₆-perfluoroalkyl,~~

~~7) —C₁-C₆-aralkyl,~~

~~8) —C₃-C₈-cycloalkyl, and~~

~~9) —heterocyclyl,~~

said alkyl, ~~aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R¹⁰;

R¹⁰ is independently selected from:

- 1) (C=O)_aO_bC₁-C₁₀ alkyl,
- 2) (C=O)_aO_baryl,
- 3) C₂-C₁₀ alkenyl,
- 4) C₂-C₁₀ alkynyl,
- 5) (C=O)_aO_b heterocyclyl,
- 6) CO₂H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O_bC₁-C₆ perfluoroalkyl,
- 11) O_a(C=O)_bNR¹²R¹³,
- 12) S(O)_mR^a,
- 13) S(O)₂NR¹²R¹³,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R¹²R¹³, or
- 17) (C=O)_aO_bC₃-C₈ cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R¹¹;

R^{10'} is halogen;

R¹¹ is selected from:

- 1) (C=O)_rO_s(C₁-C₁₀)alkyl,
- 2) O_r(C₁-C₃)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C₂-C₁₀)alkenyl,
- 8) (C₂-C₁₀)alkynyl,
- 9) (C=O)_rO_s(C₃-C₆)cycloalkyl,
- 10) (C=O)_rO_s(C₀-C₆)alkylene-aryl,

- 11) $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-heterocyclyl}$,
- 12) $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-N}(\text{R}^b)_2$,
- 13) $\text{C}(\text{O})\text{R}^a$,
- 14) $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{R}^a$,
- 15) $\text{C}(\text{O})\text{H}$,
- 16) $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{H}$, and
- 17) $\text{C}(\text{O})\text{N}(\text{R}^b)_2$,
- 18) $\text{S}(\text{O})_m\text{R}^a$, and
- 19) $\text{S}(\text{O})_2\text{N}(\text{R}^b)_2$;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from R^b , OH, $(\text{C}_1\text{-C}_6)\text{alkoxy}$, halogen, CO_2H , CN, $\text{O}(\text{C}=\text{O})\text{C}_1\text{-C}_6$ alkyl, oxo, and $\text{N}(\text{R}^b)_2$;

R^{12} and R^{13} are independently selected from:

- 1) H,
- 2) $(\text{C}=\text{O})\text{O}_b\text{C}_1\text{-C}_{10}$ alkyl,
- 3) $(\text{C}=\text{O})\text{O}_b\text{C}_3\text{-C}_8$ cycloalkyl,
- 4) $(\text{C}=\text{O})\text{O}_b\text{aryl}$,
- 5) $(\text{C}=\text{O})\text{O}_b\text{heterocyclyl}$,
- 6) $\text{C}_1\text{-C}_{10}$ alkyl,
- 7) aryl,
- 8) $\text{C}_2\text{-C}_{10}$ alkenyl,
- 9) $\text{C}_2\text{-C}_{10}$ alkynyl,
- 10) heterocyclyl,
- 11) $\text{C}_3\text{-C}_8$ cycloalkyl,
- 12) SO_2R^a , and
- 13) $(\text{C}=\text{O})\text{NR}^b_2$,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R^{11} , or

R^{12} and R^{13} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^a is independently selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^b is independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fR^{f'} or S(O)₂R^a, optionally substituted with one, two or three substituents selected from R¹¹;

R^c and R^{c'} are independently selected from: H, (C₁-C₆)alkyl, ~~aryl~~, heterocyclyl and (C₃-C₆)cycloalkyl, optionally substituted with one, two or three substituents selected from R¹¹; or

R^c and R^{c'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^d and R^{d'} are independently selected from: (C₁-C₆)alkyl, (C₁-C₆)alkoxy and NR^b₂, or

R^d and R^{d'} can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 3-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR^e, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^e is selected from: H and (C₁-C₆)alkyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^f and R^{f'} are independently selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂, or

R^f and R^{f'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R^{11} ;

R^8 is selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂.

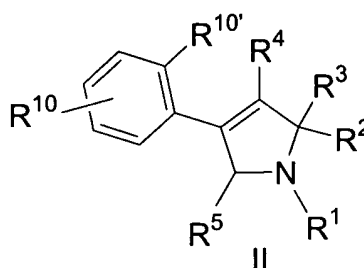
4. (Cancelled)

5. (Canceled)

6 (Original) The compound according to Claim 3 of the Formula II, or a pharmaceutically acceptable salt or stereoisomer thereof, wherein R^2 is phenyl, optionally substituted with one or two substituents selected from R^{10} .

7. (Currently amended) The compound according to Claim 1 of the formula

II:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;

b is 0 or 1;

m is 0, 1, or 2;

r is 0 or 1;

s is 0 or 1;

R^1 is selected from:

1) (C=O)C₁-C₁₀ alkyl, and

2) ~~(C=O)aryl;~~

- 3) —(C=O)C₂-C₁₀-alkenyl,
- 4) —(C=O)C₂-C₁₀-alkynyl,
- 5) —(C=O)C₃-C₈-cycloalkyl,
- 6) —(C=O)heterocyclyl,
- 2) 7) (C=O)NR^cR^{c'},
- 7) —SO₂NR^eR^{e'},
- 8) —SO₂C₁-C₁₀-alkyl,
- 9) —SO₂-aryl,
- 10) —SO₂-heterocyclyl,
- 11) —SO₂-C₃-C₈-cycloalkyl, and
- 12) —P(=O)R^dR^{d'},

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R² is phenyl, optionally substituted with one or more substituents selected from R¹⁰;

R³ is selected from:

- 1) C₁-C₁₀ alkyl-O-R^g,
- 2) C₁-C₁₀ alkyl-NR^fR^{f'},

said alkyl is optionally substituted with one or more substituents selected from R¹⁰;

R⁴ and R⁵ are independently selected from:

- 1) H, and
- 2) C₁-C₁₀ alkyl,
- 3) —aryl,
- 4) —C₂-C₁₀-alkenyl,
- 5) —C₂-C₁₀-alkynyl,
- 6) —C₁-C₆-perfluoroalkyl,
- 7) —C₁-C₆-aralkyl,
- 8) —C₃-C₈-cycloalkyl, and
- 9) —heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R¹⁰;

R¹⁰ is independently selected from:

- 1) (C=O)_aO_bC₁-C₁₀ alkyl,
- 2) (C=O)_aO_baryl,
- 3) C₂-C₁₀ alkenyl,
- 4) C₂-C₁₀ alkynyl,
- 5) (C=O)_aO_b heterocyclyl,
- 6) CO₂H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O_bC₁-C₆ perfluoroalkyl,
- 11) O_a(C=O)_bNR¹²R¹³,
- 12) S(O)_mR^a,
- 13) S(O)₂NR¹²R¹³,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R¹²R¹³, or
- 17) (C=O)_aO_bC₃-C₈ cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R¹¹;

R^{10'} is halogen;

R¹¹ is selected from:

- 1) (C=O)_rO_s(C₁-C₁₀)alkyl,
- 2) O_r(C₁-C₃)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C₂-C₁₀)alkenyl,
- 8) (C₂-C₁₀)alkynyl,
- 9) (C=O)_rO_s(C₃-C₆)cycloalkyl,
- 10) (C=O)_rO_s(C₀-C₆)alkylene-aryl,

- 11) $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-heterocyclyl}$,
- 12) $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-N}(\text{R}^b)_2$,
- 13) $\text{C}(\text{O})\text{R}^a$,
- 14) $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{R}^a$,
- 15) $\text{C}(\text{O})\text{H}$,
- 16) $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{H}$, and
- 17) $\text{C}(\text{O})\text{N}(\text{R}^b)_2$,
- 18) $\text{S}(\text{O})_m\text{R}^a$, and
- 19) $\text{S}(\text{O})_2\text{N}(\text{R}^b)_2$;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from R^b , OH, $(\text{C}_1\text{-C}_6)\text{alkoxy}$, halogen, CO_2H , CN, $\text{O}(\text{C}=\text{O})\text{C}_1\text{-C}_6$ alkyl, oxo, and $\text{N}(\text{R}^b)_2$;

R^{12} and R^{13} are independently selected from:

- 1) H,
- 2) $(\text{C}=\text{O})\text{O}_b\text{C}_1\text{-C}_{10}$ alkyl,
- 3) $(\text{C}=\text{O})\text{O}_b\text{C}_3\text{-C}_8$ cycloalkyl,
- 4) $(\text{C}=\text{O})\text{O}_b\text{aryl}$,
- 5) $(\text{C}=\text{O})\text{O}_b\text{heterocyclyl}$,
- 6) $\text{C}_1\text{-C}_{10}$ alkyl,
- 7) aryl,
- 8) $\text{C}_2\text{-C}_{10}$ alkenyl,
- 9) $\text{C}_2\text{-C}_{10}$ alkynyl,
- 10) heterocyclyl,
- 11) $\text{C}_3\text{-C}_8$ cycloalkyl,
- 12) SO_2R^a , and
- 13) $(\text{C}=\text{O})\text{NR}^b_2$,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R^{11} , or

R^{12} and R^{13} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^a is independently selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^b is independently selected from: H, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)OC₁-C₆ alkyl, (C=O)C₁-C₆ alkyl, (C=O)aryl, (C=O)heterocyclyl, (C=O)NR^fR^{f'} or S(O)₂R^a, optionally substituted with one, two or three substituents selected from R¹¹;

R^c and R^{c'} are independently selected from: H, (C₁-C₆)alkyl, ~~aryl~~, heterocyclyl and (C₃-C₆)cycloalkyl, optionally substituted with one, two or three substituents selected from R¹¹; or

R^c and R^{c'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^d and R^{d'} are independently selected from: (C₁-C₆)alkyl, (C₁-C₆)alkoxy and NR^b₂, or

R^d and R^{d'} can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 3-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR^e, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R^e is selected from: H and (C₁-C₆)alkyl, optionally substituted with one, two or three substituents selected from R¹¹;

R^f and R^{f'} are independently selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂, or

R^f and R^{f'} can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in

addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R¹¹;

R⁸ is selected from: H, (C₁-C₆)alkyl, -(C₁-C₆)alkyl-OH, -(C₁-C₆)alkyl-O-(C₁-C₆)alkyl and -(C₁-C₆)alkyl-N(R^b)₂.

8. (Amended) A compound selected from:

4-(2,5-Difluorophenyl)-2-(hydroxymethyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-(hydroxymethyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-(methoxymethyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-[(2-hydroxyethoxy)methyl]-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[(2-Aminoethoxy)methyl]-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-([2-(dimethylamino)ethyl]amino)methyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

3-{4-(2,5-Difluorophenyl)-1-[(dimethylamino)carbonyl]-2-phenyl-2,5-dihydro-1H-pyrrol-2-yl}prop-2-en-1-~~aminium~~ amine;

2-(3-Hydroxypropyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-(3-Aminopropyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-[3-(dimethylamino)propyl]-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(1-hydroxyethyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

[4-(2,5-difluorophenyl)-2-phenyl-1-(piperidin-1-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]methanol;

2-({[tert-butyl(dimethyl)silyl]oxy} methyl)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-[1-(N,N-dimethylglycyl)piperidin-4-yl]-2-(hydroxymethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(hydroxymethyl)-N-methyl-N-[1-(morpholin-4-ylacetyl)piperidin-4-yl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(hydroxymethyl)-N-methyl-2-phenyl-N-piperidin-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

N-[1-(2,2-difluoroethyl)piperidin-4-yl]-4-(2,5-difluorophenyl)-2-(hydroxymethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-[1-(2-hydroxyethyl)piperidin-4-yl]-2-(hydroxymethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-[1-(2-fluoroethyl)piperidin-4-yl]-2-(hydroxymethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(hydroxymethyl)-N-methyl-N-{1-[(methylsulfonyl)methyl]piperidin-4-yl}-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-{1-[2-fluoro-1-(fluoromethyl)ethyl]piperidin-4-yl}-2-(hydroxymethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2*S*)-*N*-(1-cyclopropylpiperidin-4-yl)-4-(2,5-difluorophenyl)-2-(hydroxymethyl)-*N*-methyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

benzyl {4-[{4-(2,5-difluorophenyl)-2-(hydroxymethyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}(methyl)amino]piperidin-1-yl} acetate;

{4-[{4-(2,5-difluorophenyl)-2-(hydroxymethyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}(methyl)amino]piperidin-1-yl} acetic acid;

methyl {4-[{4-(2,5-difluorophenyl)-2-(hydroxymethyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}(methyl)amino]piperidin-1-yl} acetate;

4-(2,5-difluorophenyl)-2-(methoxymethyl)-*N*-methyl-*N*-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxypropyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-{3-[(2,2-difluoroethyl)amino]propyl}-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-{3-[(2,2-difluoroethyl)(methyl)amino]propyl}-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-(3-aminopropyl)-4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-[3-(acetylamino)propyl]-4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-{3-[(methylsulfonyl)amino]propyl}-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

methyl 3-{4-(5-chloro-2-fluorophenyl)-1-[(dimethylamino)carbonyl]-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl} propylcarbamate;

2-{3-[(aminocarbonyl)amino]propyl}-4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

3-{4-(2,5-difluorophenyl)-1-[(dimethylamino)carbonyl]-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl}propanoic acid;

2-(3-anilino-3-oxopropyl)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydrazino-3-oxopropyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-[3-(hydroxyamino)-3-oxopropyl]-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-(2,2-difluoro-3-hydroxypropyl)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-(3-amino-2,2-difluoropropyl)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-[3-(dimethylamino)propyl]-*N*-methyl-2-phenyl-*N*-tetrahydro-2*H*-pyran-4-yl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

1-{4-(2,5-difluorophenyl)-2-[3-(dimethylamino)propyl]-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl}-2-methyl-1-oxopropan-2-ol;

3-[(2*S*)-1-[(2*S*)-2-amino-2-cyclopropylethanoyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl]-*N,N*-dimethylpropan-1-amine; and

(2*S*)-2-(3-amino-4,4-difluorobutyl)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

or a pharmaceutically acceptable salt or stereoisomer thereof.

9. (Original) The compound according to Claim 8 which is selected from:

4-(2,5-Difluorophenyl)-2-(hydroxymethyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[(2-Aminoethoxy)methyl]-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-(3-Aminopropyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-2-[3-(dimethylamino)propyl]-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

or a pharmaceutically acceptable salt or stereoisomer thereof.

10. (Canceled)

11. (Original) The compound according to Claim 1 selected from:

2-[(2-Aminoethoxy)methyl]-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide hydrochloride salt;

4-(2,5-Difluorophenyl)-2-([2-(dimethylamino)ethyl]amino)methyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide bis TFA salt;

2-(3-Aminopropyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide hydrochloride salt;

4-(2,5-Difluorophenyl)-2-[3-(dimethylamino)propyl]-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide TFA salt;

4-(2,5-difluorophenyl)-N-[1-(glycyl)piperidin-4-yl]-2-(hydroxymethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide TFA salt;

3-[(2*S*)-1-[(2*S*)-2-amino-2-cyclopropylethanoyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl]-*N,N*-dimethylpropan-1-amine bis-TFA salt;

3-[(2*R*)-1-[(2*S*)-2-amino-2-cyclopropylethanoyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-2-yl]-*N,N*-dimethylpropan-1-amine bis-TFA salt;

4-(2,5-difluorophenyl)-2-[3-(dimethylamino)propyl]-*N*-methyl-*N*-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide bis-TFA salt;

4-(2,5-difluorophenyl)-2-[3-(ethylamino)propyl]-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide TFA salt;

4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2-{3-[(pyridin-4-ylmethyl)amino]propyl}-2,5-dihydro-1*H*-pyrrole-1-carboxamide bis-TFA salt; and

4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-(3-[[4-methyl-1*H*-imidazol-2-yl)methyl]amino}propyl)-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide bis-TFA salt.

12. (Original) A pharmaceutical composition that is comprised of a compound in accordance with Claim 1 and a pharmaceutically acceptable carrier.

13. (Cancelled)

14. (Cancelled)

15. (Cancelled)

16. (Canceled)

17. (Canceled)

18. (Canceled)

19. (Canceled)
20. (Canceled)
21. (Canceled)
22. (Canceled)
23. (Canceled)
24. (Canceled)
25. (Cancelled)
26. (Cancelled)
27. (Canceled)
28. (Canceled)
29. (Canceled)
30. (Canceled)
31. (Canceled)
32. (Canceled)
33. (Canceled)
34. (Canceled)
35. (Canceled)

36. (Canceled)

37. (Canceled)

38. (Cancelled)